

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082015\
 Data File : BF081130.D
 Acq On : 20 Aug 2015 18:06
 Operator : UM/IZ
 Sample : PB85099BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85099BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.699	243	246	248	rBV	687568	978135	51.70%	6.821%
2	5.144	283	285	288	rBV	1074436	1069413	56.52%	7.458%
3	6.219	376	379	381	rBV	1309570	1249289	66.03%	8.712%
4	6.333	387	389	392	rBV	1063790	1035633	54.73%	7.222%
5	6.573	407	410	412	rBV	516740	436822	23.09%	3.046%
6	6.733	421	424	426	rBV	1042751	1230986	65.06%	8.585%
7	7.145	457	460	462	rBV	1144166	1158719	61.24%	8.081%
8	7.865	520	523	525	rBV	407606	560268	29.61%	3.907%
9	8.939	614	617	620	rBV	1745612	1706798	90.21%	11.903%
10	9.602	673	675	678	rBB	572631	593189	31.35%	4.137%
11	10.391	741	744	746	rBV	698448	725624	38.35%	5.060%
12	11.076	801	804	806	rBV	798696	672945	35.57%	4.693%
13	12.665	940	943	945	rBV	1943643	1892094	100.00%	13.195%
14	13.694	1030	1033	1035	rBV	683867	613510	32.42%	4.278%
15	15.031	1147	1150	1153	rBV	390349	415994	21.99%	2.901%

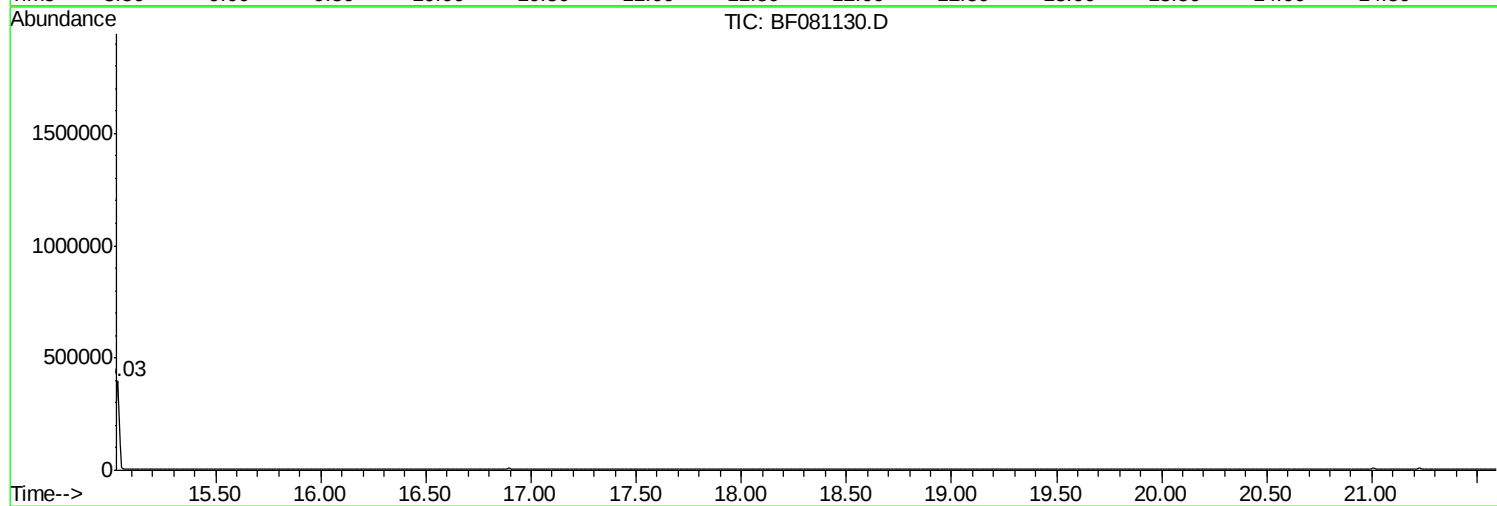
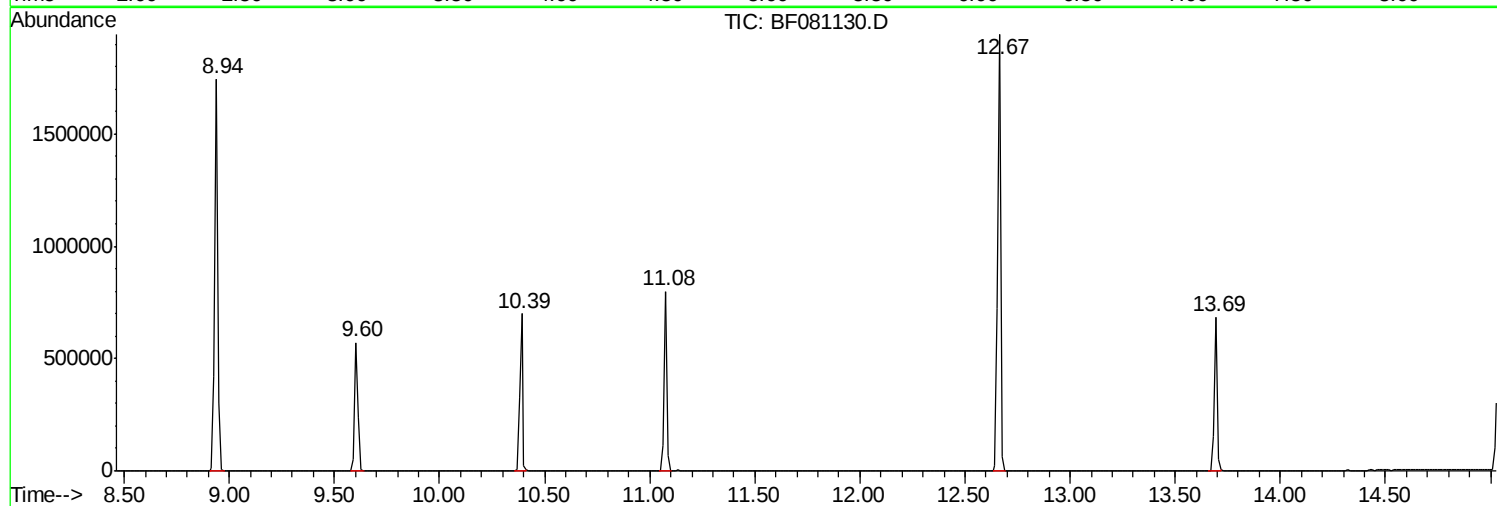
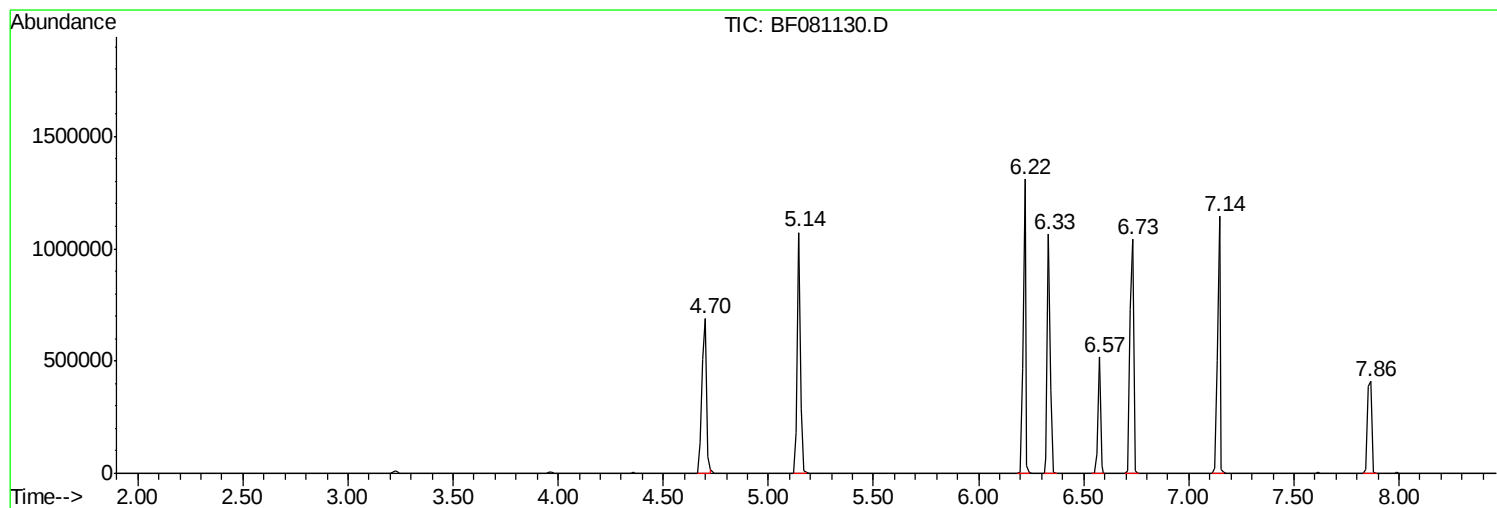
Sum of corrected areas: 14339419

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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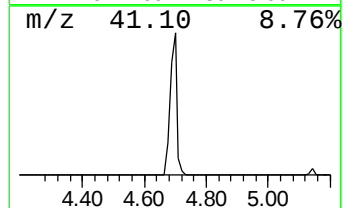
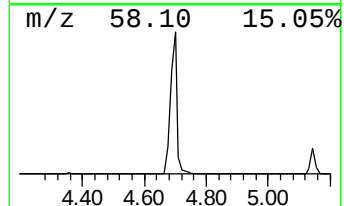
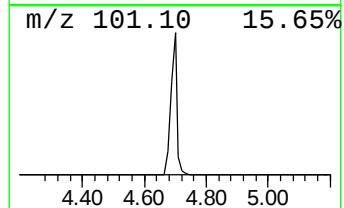
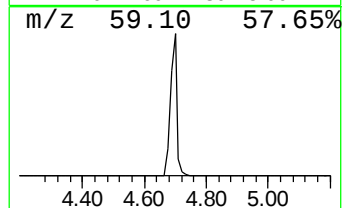
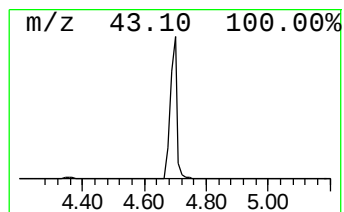
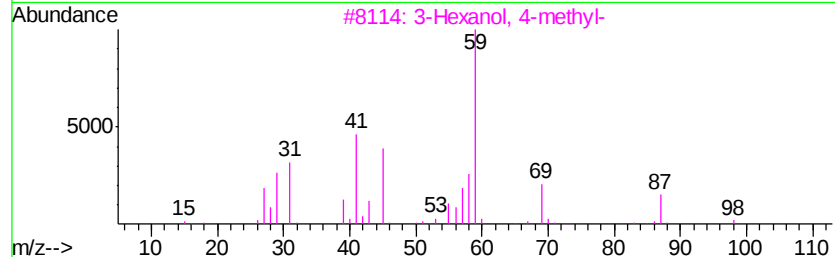
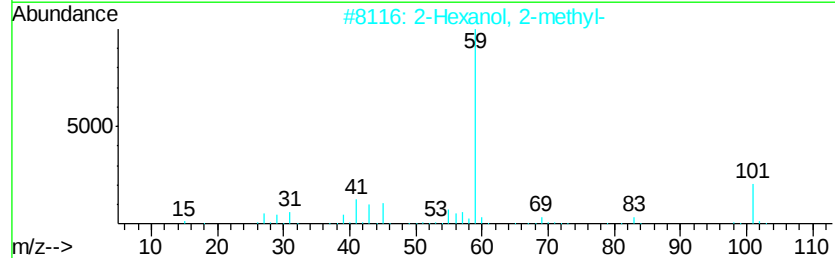
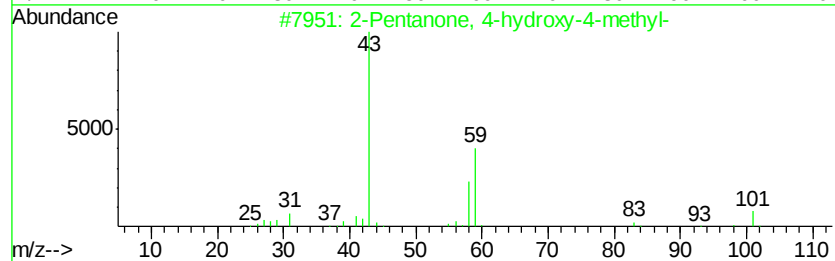
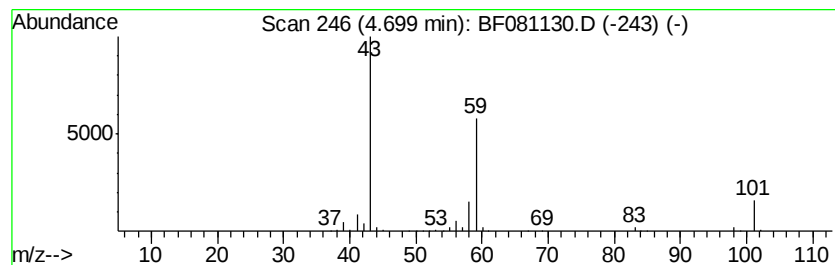
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.70	44.78 ng	978135	1,4-Dichlorobenzene-d4	6.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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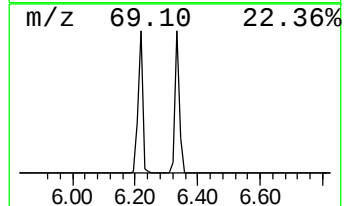
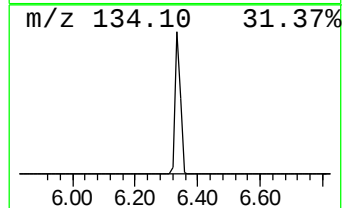
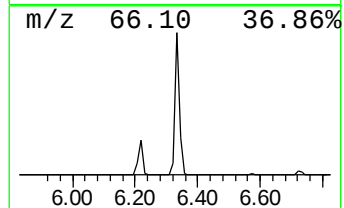
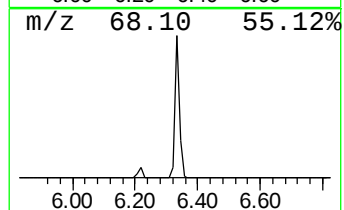
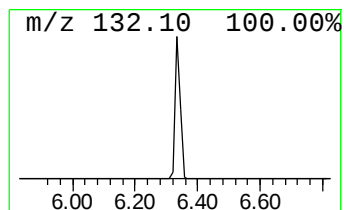
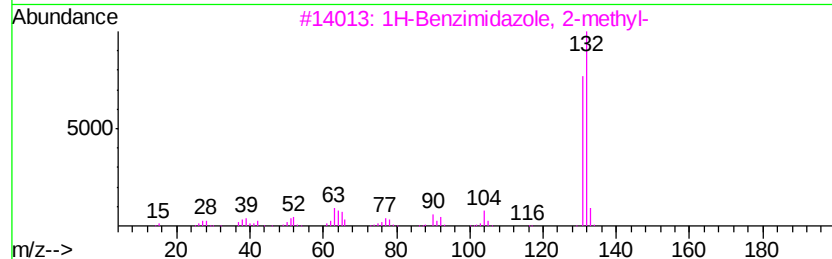
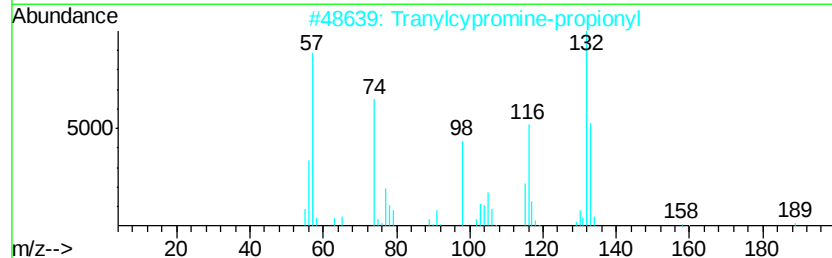
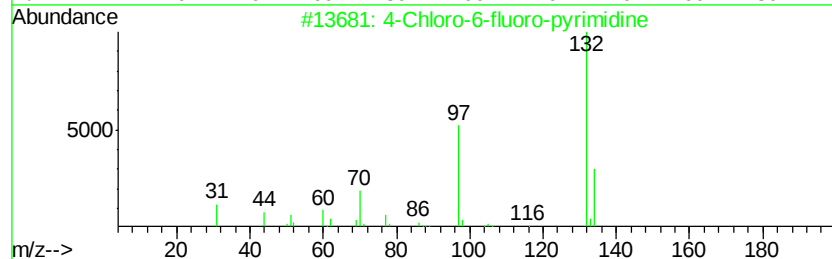
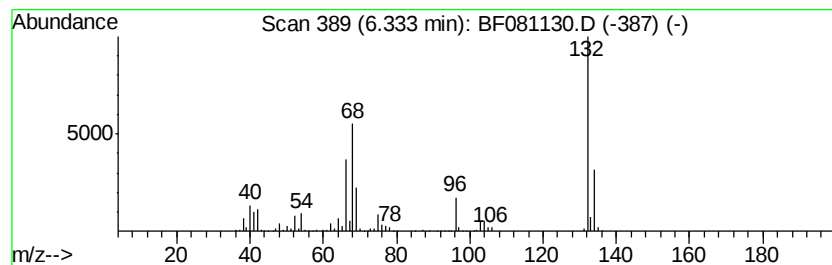
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Peak Number 2 unknown6.33 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	47.42 ng	1035630	1,4-Dichlorobenzene-d4	6.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.70	44.8	ng	978135	1	6.57	436822	20.0
unknown6.33	6.33	47.4	ng	1035630	1	6.57	436822	20.0